Propanamide, 3-phenyl-N-ethyl-N-tetradecyl-

Inchi: InChI=1S/C25H43NO/c1-3-5-6-7-8-9-10-11-12-13-14-18-23-26(4-2)25(27)22-21-24-19-1
InchiKey: YZHRGGVCMRKQDV-UHFFFAOYSA-N
Formula: C25H43NO
SMILES: CCCCCCCCCCCCCCN(CC)C(=O)CCc1ccccc1
Mol. weight [g/mol]: 373.62

Physical Properties

Property code	Value	Unit	Source
gf	253.89	kJ/mol	Joback Method
hf	-367.85	kJ/mol	Joback Method
hfus	59.17	kJ/mol	Joback Method
hvap	82.31	kJ/mol	Joback Method
log10ws	-7.74		Crippen Method
logp	7.169		Crippen Method
mcvol	350.900	ml/mol	McGowan Method
pc	960.88	kPa	Joback Method
rinpol	3320.00		NIST Webbook
tb	864.39	K	Joback Method
tc	1060.51	K	Joback Method
tf	480.33	K	Joback Method
vc	1.351	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	1141.72	J/mol×K	864.39	Joback Method
cpg	1161.77	J/mol×K	897.08	Joback Method
cpg	1180.66	J/mol×K	929.76	Joback Method
cpg	1198.45	J/mol×K	962.45	Joback Method
cpg	1215.22	J/mol×K	995.14	Joback Method
cpg	1231.03	J/mol×K	1027.83	Joback Method
cpg	1245.96	J/mol×K	1060.51	Joback Method

Sources

Crippen Method:	http://pubs.acs.org/doi/abs/10.1021/ci990307l
Crippen Method:	https://www.chemeo.com/doc/models/crippen_log10ws
Joback Method:	https://en.wikipedia.org/wiki/Joback_method
McGowan Method:	http://link.springer.com/article/10.1007/BF02311772
NIST Webbook:	http://webbook.nist.gov/cgi/cbook.cgi?ID=U415404&Units=SI

Legend

cpg:	Ideal gas heat capacity
gf:	Standard Gibbs free energy of formation
hf:	Enthalpy of formation at standard conditions
hfus:	Enthalpy of fusion at standard conditions
hvap:	Enthalpy of vaporization at standard conditions
log10ws:	Log10 of Water solubility in mol/l
logp:	Octanol/Water partition coefficient
mcvol:	McGowan's characteristic volume
pc:	Critical Pressure
rinpol:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume